

Data Path : Z:\svoasrv\HPCHEM1\BNA_F\Data\BF032124\
 Data File : BF137675.D
 Acq On : 21 Mar 2024 22:23
 Operator : CG\JU
 Sample : P1804-01DL 10X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 JC-03-03192024-ADL

Quant Time: Mar 22 01:45:28 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_F\Methods\8270-BF022924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 14 14:45:20 2024
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By :Jagrut Upadhyay 03/22/2024
 Supervised By :mohammad ahmed 03/26/2024

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	6.728	152	114198	20.000	ng	0.00
21) Naphthalene-d8	8.010	136	416504	20.000	ng	0.00
39) Acenaphthene-d10	9.757	164	199339	20.000	ng	0.00
64) Phenanthrene-d10	11.239	188	322356	20.000	ng	0.00
76) Chrysene-d12	13.880	240	274140	20.000	ng	0.00
86) Perylene-d12	15.304	264	206955	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.334	112	48389	6.418	ng	0.00
7) Phenol-d6	6.369	99	59307	5.449	ng	0.00
23) Nitrobenzene-d5	7.298	82	30312	3.382	ng	0.00
42) 2,4,6-Tribromophenol	10.551	330	10619	6.066	ng	0.00
45) 2-Fluorobiphenyl	9.081	172	80962	6.358	ng	0.00
79) Terphenyl-d14	12.827	244	92008	4.615	ng	0.00
Target Compounds						Qvalue
31) Naphthalene	8.028	128	80308	3.594	ng	98
37) 2-Methylnaphthalene	8.722	142	28834	2.024	ng	99
52) Acenaphthene	9.792	154	40964	3.854	ng	98
55) Dibenzofuran	9.969	168	39816	2.460	ng	97
58) Fluorene	10.304	166	61942	4.983	ng	99
71) Phenanthrene	11.269	178	465584	26.539	ng	98
72) Anthracene	11.316	178	157618	8.748	ng	97
73) Carbazole	11.480	167	39168	2.538	ng	100
75) Fluoranthene	12.451	202	695166	40.539	ng	99
78) Pyrene	12.680	202	617777	22.966	ng	99
81) Benzo(a)anthracene	13.868	228	317063	16.501	ng	99
83) Chrysene	13.904	228	253754	13.067	ng	96
87) Indeno(1,2,3-cd)pyrene	16.692	276	94887	6.967	ng	94
88) Benzo(b)fluoranthene	14.898	252	271130m	22.147	ng	
89) Benzo(k)fluoranthene	14.921	252	74907m	5.697	ng	
90) Benzo(a)pyrene	15.245	252	183452	15.206	ng	98
91) Dibenzo(a,h)anthracene	16.698	278	22235	2.013	ng #	91
92) Benzo(g,h,i)perylene	17.109	276	93496	8.240	ng	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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